L Number	Hits	Search Text	DB	Time stamp
5	424	562/439.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05
6	190	548/480.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05
7	256	549/69.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05
8	246	548/473.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05
10	903	548/253.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05
11	276	549/373.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05
12	2244	562/439.ccls. or 548/480.ccls. or 549/69.ccls. or 548/473.ccls. or 588/417.ccls. or 548/253.ccls. or 549/373.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05
13	614	514/452.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05
14	1002	514/381.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05
15	998	514/563.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05
16	85	514/113.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05
17	233	514/447.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05
18	258	514/417.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05
19	654	514/364.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05
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22	6894	glucagon	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05

1	10	US-5880139-$.DID. OR US-5776954-$.DID. OR US-4359474-$.DID. OR US-4374130-$.DID. OR US-5837719-$.DID.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05
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2	2	6096773.pn.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05
3	2	6262084.pn.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05
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27	2	6503949.pn.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 08:06

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3	BRS	L7	256	549/69.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05		
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6	BRS	L11	276	549/373.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05		
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9	BRS	L14	1002	514/381.ccls.	USPAT; US-PG PUB; EPO; JPO; DERWENT	2004/07/16 06:05		
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18	BRS	L1	10	US-5880139-$.DID. OR US-5776954-$.DID. OR US-4359474-$.DID. OR US-4374130-$.DID. OR US-5837719-$.DID.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05		
19	BRS	L24	91	((514/452.ccls. or 514/381.ccls. or 514/563.ccls. or 514/113.ccls. or 514/447.ccls. or 514/417.ccls. or 514/364.ccls.) or (562/439.ccls. or 548/480.ccls. or 549/69.ccls. or 548/473.ccls. or 588/417.ccls. or 548/253.ccls. or 549/373.ccls.)) and glucagon	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/07/16 06:05		
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23	BRS	L23	477	(514/452.ccls. or 514/381.ccls. or 514/563.ccls. or 514/113.ccls. or 514/447.ccls. or 514/417.ccls. or 514/364.ccls.) and (562/439.ccls. or 548/480.ccls. or 549/69.ccls. or 548/473.ccls. or 588/417.ccls. or 548/253.ccls. or 549/373.ccls.)	USPAT; US-PG PUB; EPO; JPO; DERWENT	2004/07/16 06:05		
24	BRS	L25	2	6562807.pn.	USPAT; US-PG PUB; EPO; JPO; DERWENT	2004/07/16 07:59		
25	BRS	L26	1	6762318.pn.	USPAT; US-PG PUB; EPO; JPO; DERWENT	2004/07/16 08:06		
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=> file reg
COST IN U.S. DOLLARS

	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.21	0.21

FILE 'REGISTRY' ENTERED AT 06:14:46 ON 16 JUL 2004
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

Property values tagged with IC are from the ZIC/VINITI data file
provided by InfoChem.

STRUCTURE FILE UPDATES: 14 JUL 2004 HIGHEST RN 710278-45-8
DICTIONARY FILE UPDATES: 14 JUL 2004 HIGHEST RN 710278-45-8

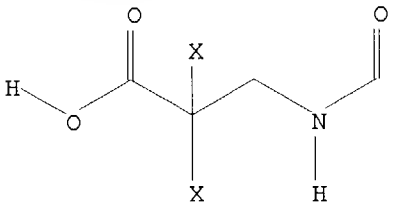
TSCA INFORMATION NOW CURRENT THROUGH JANUARY 6, 2004

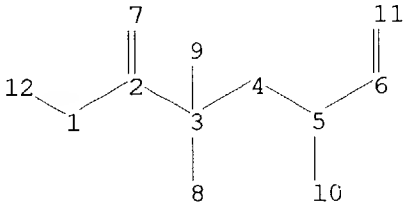
Please note that search-term pricing does apply when
conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more
information enter HELP PROP at an arrow prompt in the file or refer
to the file summary sheet on the web at:
http://www.cas.org/ONLINE/DBSS/registryss.html

=>
Uploading C:\Examination Auxillary files\09995987\09995987 third try fixed H broader.str





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chain bonds :
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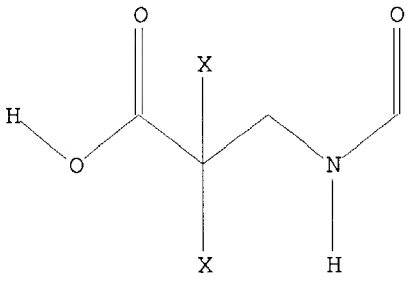
exact/norm bonds :
4-5 5-6 6-11
exact bonds :
1-12 2-3 3-4 3-8 3-9 5-10
normalized bonds :
1-2 2-7

G1:C,O,S,N

Hydrogen count :
4:>= minimum 2
Match level :
1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS 8:CLASS 9:CLASS
10:CLASS 11:CLASS 12:CLASS

L1 STRUCTURE UPLOADED

=> d l1
L1 HAS NO ANSWERS
L1 STR



G1 C,O,S,N

Structure attributes must be viewed using STN Express query preparation.

=> search l1 sss sam
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SAMPLE SCREEN SEARCH COMPLETED - 21 TO ITERATE

100.0% PROCESSED 21 ITERATIONS 1 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 146 TO 694
PROJECTED ANSWERS: 1 TO 80

L2 1 SEA SSS SAM L1

=> d scan

L2 1 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Propanoic acid, 3-[[4-[[[[(3,5-dichlorophenyl)amino]carbonyl][4-[4-(1,1-dimethylethyl)-1-cyclohexen-1-yl]phenyl]amino]methyl]benzoyl]amino]-2,2-difluoro- (9CI)
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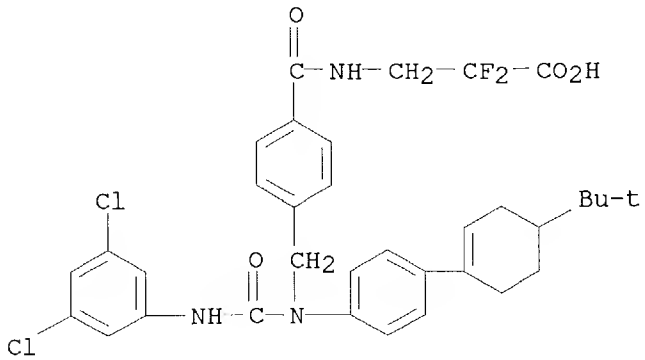

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> search l1 sss full
FULL SEARCH INITIATED 06:15:46 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 525 TO ITERATE

100.0% PROCESSED 525 ITERATIONS 3 ANSWERS
SEARCH TIME: 00.00.01

L3 3 SEA SSS FUL L1

=> d scan

L3 3 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Propanoic acid, 3-[[4-[[[[(3,5-dichlorophenyl)amino]carbonyl][4-[4-(1,1-dimethylethyl)-1-cyclohexen-1-yl]phenyl]amino]methyl]benzoyl]amino]-2,2-difluoro- (9CI)
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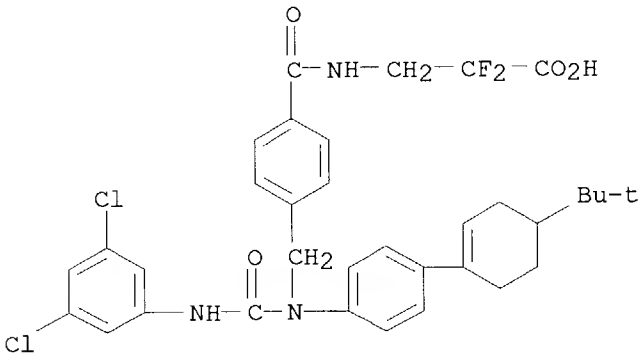

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):3

L3 3 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Propanoic acid, 2,2-difluoro-3-[[(phenylmethoxy)carbonyl]amino]- (9CI)
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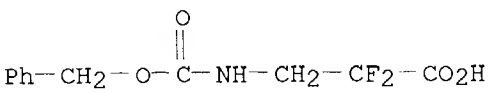
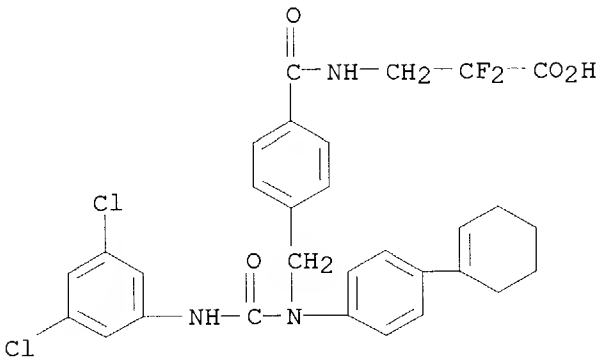


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 3 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Propanoic acid, 3-[[4-[[[4-(1-cyclohexen-1-yl)phenyl][[(3,5-

dichlorophenyl)amino]carbonyl]amino]methyl]benzoyl]amino]-2,2-difluoro-(9CI)
MF C30 H27 Cl2 F2 N3 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	155.84	156.05

FILE 'CAPLUS' ENTERED AT 06:16:02 ON 16 JUL 2004
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FILE COVERS 1907 - 16 Jul 2004 VOL 141 ISS 4
FILE LAST UPDATED: 15 Jul 2004 (20040715/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> l3
L4 1 L3

=> d l4 ti fbib abs

L4 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2004 ACS on STN
TI Preparation of ureidomethylbenzoylaminodifluoropropionates as glucagon antagonists/inverse agonists.

AN 2002:391685 CAPLUS
DN 136:385945
TI Preparation of ureidomethylbenzoylaminodifluoropropionates as glucagon antagonists/inverse agonists.
IN Jorgensen, Anker Steen; Madsen, Peter
PA Novo Nordisk A/S, Den.
SO PCT Int. Appl., 85 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2002040446	A1	20020523	WO 2001-DK760	20011115
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	RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
				DK 2000-1733	A 20001117
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				DK 2000-1733	A 20001117
				WO 2001-DK760	W 20011115
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	US 2003027849	A1	20030206	US 2001-995987	20011116
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				US 2000-252322PP	20001120

OS MARPAT 136:385945
AB HO2CCF2CH2NHCOZCHR2N(E)XD [R2 = H, alkyl; Z = (substituted) arylene, heteroarylene; X = (CH2)q(CR12R13)r(CH2)s, CO(CR12R13)r(CH2)s, NR11CO(CR12R13)r(CH2)s, etc.; r = 0, 1; s = 0-3; R11, R12, R13 = H , alkyl; D = (substituted) Ph, naphthyl, pyridyl, indenyl, benzothienyl, thienyl, furyl, benzofuryl, etc.; E = (substituted) cyclohexyl, Ph, PhCH2, PhCH2CH2, indanyl, benzhydryl, etc.], were prepared Thus, Me 4-[(4-cyclohex-1-enylphenylamino)methyl]benzoate (preparation given) in CH2Cl2 containing diisopropylethylamine was treated with 3,5-dichlorophenyl isocyanate to give a residue which was saponified with LiOH. The resulting acid in DMF was treated with 3-[(dimethyliminium)(dimethylamino)methyl]-1,2,3-benzotriazol-1-ium-1-olate hexafluorophosphate, diisopropylethylamine, Me 3-amino-2,2-difluoropropionate hydrochloride to give the uncharacterized amide ester, which was saponified with aqueous LiOH in THF/MeOH to give 3-[4-[1-(4-cyclohex-1-enylphenyl)-3-(3,5-dichlorophenyl)ureidomethyl]benzoylamino]-2,2-difluoropropionic acid. In a human glucagon receptor binding assay, title compds. showed IC50<1000 nM.
RE.CNT 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> logoff hold
COST IN U.S. DOLLARS SINCE FILE ENTRY TOTAL SESSION

FULL ESTIMATED COST	3.14	159.19

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-0.74	-0.74

SESSION WILL BE HELD FOR 60 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 06:16:31 ON 16 JUL 2004

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LOGINID:SSSPTA1623PAZ

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* * * * * * * * * * Welcome to STN International * * * * * * * * * *

NEWS	1		Web Page URLs for STN Seminar Schedule - N. America
NEWS	2		"Ask CAS" for self-help around the clock
NEWS	3	May 12	EXTEND option available in structure searching
NEWS	4	May 12	Polymer links for the POLYLINK command completed in REGISTRY
NEWS	5	May 27	New UPM (Update Code Maximum) field for more efficient patent SDIs in CAplus
NEWS	6	May 27	CAplus super roles and document types searchable in REGISTRY
NEWS	7	Jun 22	STN Patent Forums to be held July 19-22, 2004
NEWS	8	Jun 28	Additional enzyme-catalyzed reactions added to CASREACT
NEWS	9	Jun 28	ANTE, AQUALINE, BIOENG, CIVILENG, ENVIROENG, MECHENG, and WATER from CSA now available on STN(R)
NEWS	10	Jul 12	BEILSTEIN enhanced with new display and select options, resulting in a closer connection to BABS

NEWS EXPRESS	MARCH 31 CURRENT WINDOWS VERSION IS V7.00A, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 26 APRIL 2004
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NEWS INTER	General Internet Information
NEWS LOGIN	Welcome Banner and News Items
NEWS PHONE	Direct Dial and Telecommunication Network Access to STN
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* * * * * * * * * * * * * STN Columbus * * * * * * * * * * * * * * * *

FILE 'HOME' ENTERED AT 07:23:23 ON 16 JUL 2004

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.21	0.21

FILE 'CAPLUS' ENTERED AT 07:23:45 ON 16 JUL 2004
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FILE COVERS 1907 - 16 Jul 2004 VOL 141 ISS 4
FILE LAST UPDATED: 15 Jul 2004 (20040715/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> file reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.46	0.67

FILE 'REGISTRY' ENTERED AT 07:23:49 ON 16 JUL 2004
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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 14 JUL 2004 HIGHEST RN 710278-45-8
DICTIONARY FILE UPDATES: 14 JUL 2004 HIGHEST RN 710278-45-8

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 6, 2004

Please note that search-term pricing does apply when conducting SmartSELECT searches.

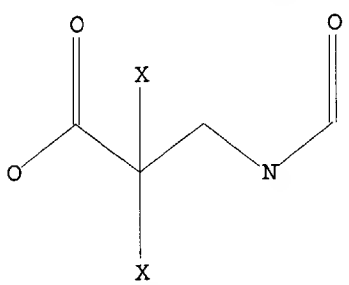
Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at:
http://www.cas.org/ONLINE/DBSS/registryss.html

=>
Uploading C:\Examination Auxillary files\09995987\09995987 patetable core broadest.str

L1 STRUCTURE UPLOADED

=> d l1
L1 HAS NO ANSWERS
L1 STR



G1 C,O,S,N

Structure attributes must be viewed using STN Express query preparation.

=> search l1 sss sam
SAMPLE SEARCH INITIATED 07:24:18 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 194 TO ITERATE

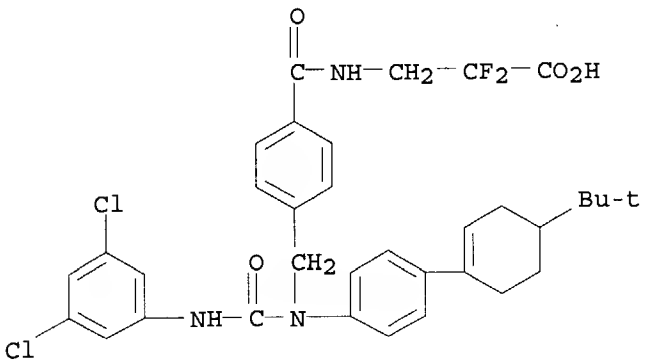
100.0% PROCESSED 194 ITERATIONS 2 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 3045 TO 4715
PROJECTED ANSWERS: 2 TO 124

L2 2 SEA SSS SAM L1

=> d scan

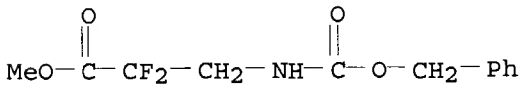
L2 2 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Propanoic acid, 3-[[4-[[[[(3,5-dichlorophenyl)amino]carbonyl][4-[4-(1,1-dimethylethyl)-1-cyclohexen-1-yl]phenyl]amino]methyl]benzoyl]amino]-2,2-difluoro- (9CI)
MF C34 H35 Cl2 F2 N3 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):2

L2 2 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Propanoic acid, 2,2-difluoro-3-[[(phenylmethoxy)carbonyl]amino]-, methyl ester (9CI)
MF C12 H13 F2 N O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

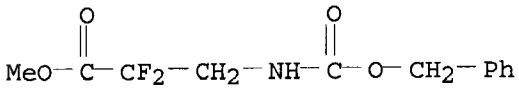
=> search l1 sss full
FULL SEARCH INITIATED 07:24:40 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 4048 TO ITERATE

100.0% PROCESSED 4048 ITERATIONS 4 ANSWERS
SEARCH TIME: 00.00.01

L3 4 SEA SSS FUL L1

=> d scan

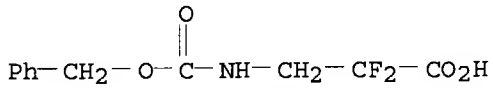
L3 4 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Propanoic acid, 2,2-difluoro-3-[[(phenylmethoxy)carbonyl]amino]-, methyl ester (9CI)
MF C12 H13 F2 N O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

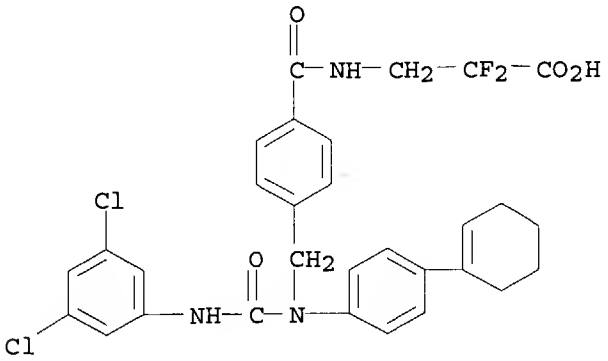
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):4

L3 4 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Propanoic acid, 2,2-difluoro-3-[[(phenylmethoxy)carbonyl]amino]- (9CI)
MF C11 H11 F2 N O4



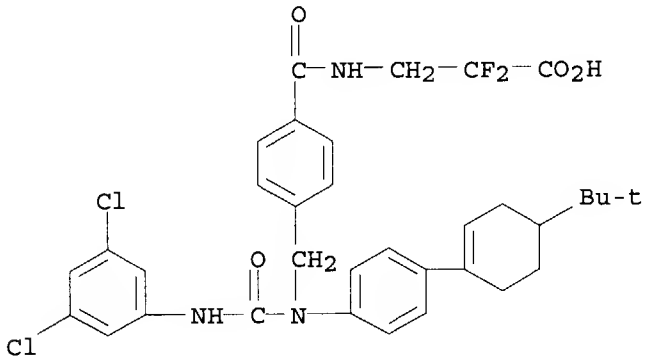
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 4 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Propanoic acid, 3-[[4-[[[4-(1-cyclohexen-1-yl)phenyl][[(3,5-dichlorophenyl)amino]carbonyl]amino]methyl]benzoyl]amino]-2,2-difluoro-(9CI)
MF C30 H27 Cl2 F2 N3 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 4 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Propanoic acid, 3-[[4-[[[[(3,5-dichlorophenyl)amino]carbonyl][4-[4-(1,1-dimethylethyl)-1-cyclohexen-1-yl]phenyl]amino]methyl]benzoyl]amino]-2,2-difluoro- (9CI)
MF C34 H35 Cl2 F2 N3 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	155.84	156.51

FILE 'CAPLUS' ENTERED AT 07:25:00 ON 16 JUL 2004
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FILE COVERS 1907 - 16 Jul 2004 VOL 141 ISS 4
FILE LAST UPDATED: 15 Jul 2004 (20040715/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> l3
L4 1 L3

=> d l4

L4 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2004 ACS on STN
AN 2002:391685 CAPLUS
DN 136:385945
TI Preparation of ureidomethylbenzoylaminodifluoropropionates as glucagon antagonists/inverse agonists.
IN Jorgensen, Anker Steen; Madsen, Peter
PA Novo Nordisk A/S, Den.
SO PCT Int. Appl., 85 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2002040446	A1	20020523	WO 2001-DK760	20011115
	W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PH, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
	RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
	AU 2002023502	A5	20020527	AU 2002-23502	20011115
	EP 1345891	A1	20030924	EP 2001-996529	20011115
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR				
	JP 2004513936	T2	20040513	JP 2002-542774	20011115
	US 2003027849	A1	20030206	US 2001-995987	20011116
PRAI	DK 2000-1733	A	20001117		
	US 2000-252322P	P	20001120		
	WO 2001-DK760	W	20011115		

OS MARPAT 136:385945
RE.CNT 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> file beilstein
COST IN U.S. DOLLARS

	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	2.03	158.54

FILE 'BEILSTEIN' ENTERED AT 07:26:00 ON 16 JUL 2004

FILE RELOADED ON OCTOBER 20, 2002
FILE LAST UPDATED ON JUNE 15, 2004

FILE COVERS 1771 TO 2003.
*** FILE CONTAINS 8,997,153 SUBSTANCES ***

>>> PLEASE NOTE: Reaction data and substance data are stored in separate documents and can not be searched together in one query.
Reaction data for BEILSTEIN compounds may be displayed immediately with the display codes PRE (preparations) and REA (reactions). A substance answer set retrieved after the search for a chemical name, a molecular formula or a structure search for example can be restricted to compounds with available reaction information by concatenation with PRE/FA, REA/FA or more general with RX/FA. The BEILSTEIN Registry Number (BRN) is the link between a BEILSTEIN compound and belonging reactions. For more detailed reaction searches BRNs can be selected from substance answer sets and searched in the next step as reaction partner BRNs - Reactant (RX.RBRN) or Product BRN (RX.PBRN). After a search for reaction details substance documents associated with reactants or products may be retrieved by searching RX.PBRNs or RX.RBRNs as BRNs. <<<

>>> FOR SEARCHING PREPARATIONS SEE HELP PRE <<<

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**********************************************************************
* PLEASE NOTE THAT THERE ARE NO FORMATS FREE OF COST.                *
* SET NOTICE FEATURE: THE COST ESTIMATES CALCULATED FOR SET NOTICE   *
* ARE BASED ON THE HIGHEST PRICE CATEGORY. THEREFORE; THESE          *
* ESTIMATES MAY NOT REFLECT THE ACTUAL COSTS.                        *
* FOR PRICE INFORMATION SEE HELP COST                                *
**********************************************************************
```

=> l1
SAMPLE SEARCH INITIATED 07:26:08 FILE 'BEILSTEIN'
SAMPLE SCREEN SEARCH COMPLETED - 67 TO ITERATE

100.0% PROCESSED 67 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.02

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 849 TO 1831
PROJECTED ANSWERS: 0 TO 0

L5 0 SEA SSS SAM L1